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LOGINID: SSPTAMLL1621
PASSWORD:
TERMINAL (ENTER 1, 2, 3, OR ?):2
                     Welcome to STN International
                 Web Page URLs for STN Seminar Schedule - N. America
NEWS
      1
                  "Ask CAS" for self-help around the clock
 NEWS
                 CA/CAplus pre-1967 chemical substance index entries enhanced
 NEWS 3
         DEC 18
                 with preparation role
         DEC 18
                 CA/CAplus patent kind codes updated
 NEWS
                 MARPAT to CA/CAplus accession number crossover limit increased
         DEC 18
 NEWS 5
                  to 50,000
                 MEDLINE updated in preparation for 2007 reload
         DEC 18
 NEWS 6
         DEC 27
                 CA/CAplus enhanced with more pre-1907 records
      7
 NEWS
         JAN 08
                 CHEMLIST enhanced with New Zealand Inventory of Chemicals
 NEWS 8
                 CA/CAplus Company Name Thesaurus enhanced and reloaded
 NEWS 9 JAN 16
NEWS 10 JAN 16
                 IPC version 2007.01 thesaurus available on STN
 NEWS 11 JAN 16
                 WPIDS/WPINDEX/WPIX enhanced with IPC 8 reclassification data
 NEWS 12 JAN 22
                 CA/CAplus updated with revised CAS roles
                  CA/CAplus enhanced with patent applications from India
 NEWS 13 JAN 22
                  PHAR reloaded with new search and display fields
 NEWS 14 JAN 29
                  CAS Registry Number crossover limit increased to 300,000 in
 NEWS 15 JAN 29
                  multiple databases
                  PATDPASPC enhanced with Drug Approval numbers
 NEWS 16 FEB 15
 NEWS 17 FEB 15
                 RUSSIAPAT enhanced with pre-1994 records
 NEWS 18 FEB 23 KOREAPAT enhanced with IPC 8 features and functionality
 NEWS 19 FEB 26 MEDLINE reloaded with enhancements
 NEWS 20 FEB 26 EMBASE enhanced with Clinical Trial Number field
 NEWS 21 FEB 26
                 TOXCENTER enhanced with reloaded MEDLINE
 NEWS 22 FEB 26
                  IFICDB/IFIPAT/IFIUDB reloaded with enhancements
 NEWS 23 FEB 26 CAS Registry Number crossover limit increased from 10,000
                  to 300,000 in multiple databases
 NEWS 24 MAR 15 WPIDS/WPIX enhanced with new FRAGHITSTR display format
 NEWS 25 MAR 16 CASREACT coverage extended
 NEWS 26 MAR 20 MARPAT now updated daily
 NEWS 27 MAR 22 LWPI reloaded
 NEWS EXPRESS NOVEMBER 10 CURRENT WINDOWS VERSION IS V8.01c, CURRENT
               MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP),
               AND CURRENT DISCOVER FILE IS DATED 25 SEPTEMBER 2006.
               STN Operating Hours Plus Help Desk Availability
 NEWS HOURS
               Welcome Banner and News Items
 NEWS LOGIN
               For general information regarding STN implementation of IPC 8
 NEWS IPC8
 NEWS X25
               X.25 communication option no longer available
```

Enter NEWS followed by the item number or name to see news on that specific topic.

Page 1 searched 9/5/07 updated str search

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=> fil reg COST IN U.S. DOLLARS

FULL ESTIMATED COST

SINCE FILE TOTAL ENTRY SESSION 0.21 0.21

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STRUCTURE FILE UPDATES: 23 MAR 2007 HIGHEST RN 928114-47-0 DICTIONARY FILE UPDATES: 23 MAR 2007 HIGHEST RN 928114-47-0

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http://www.cas.org/ONLINE/UG/regprops.html

=> s 335200-36-7/rn or 845785-97-9/rn or 845785-98-0/rn or 845785-99-1/rn or 845786-00-7/rn or 845786-01-8/rn or 845786-02-9/rn or 845786-03-0/rn or 845786-04-1/rn

1 335200-36-7/RN

1 845785-97-9/RN

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1 845785-99-1/RN

1 845786-00-7/RN

1 845786-01-8/RN

1 845786-02-9/RN

1 845786-03-0/RN

1 845786-04-1/RN

L1 9 335200-36-7/RN OR 845785-97-9/RN OR 845785-98-0/RN OR 845785-99-1/RN OR 845786-00-7/RN OR 845786-01-8/RN OR 845786-02-9/ RN OR 845786-03-0/RN OR 845786-04-1/RN

=> d scan

L1 9 ANSWERS REGISTRY COPYRIGHT 2007 ACS on STN

IN Butanedioic acid, (4-bromophenyl)-, 1-(phenylmethyl) ester (9CI)

MF C17 H15 Br O4

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):9

L1 9 ANSWERS REGISTRY COPYRIGHT 2007 ACS on STN

IN Butanedioic acid, (4-bromophenyl)-, 1-(1,1-dimethylethyl) 4-(phenylmethyl)

ester (9CI)

MF C21 H23 Br O4

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L1 9 ANSWERS REGISTRY COPYRIGHT 2007 ACS on STN

IN Butanedioic acid, (4-bromophenyl)-, 4-(1,1-dimethylethyl) 1-(phenylmethyl)

ester (9CI)

MF C21 H23 Br O4

Page 3 searched 9/5/07 updated str search

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L1 9 ANSWERS REGISTRY COPYRIGHT 2007 ACS on STN

IN Benzeneacetic acid, 4-[[(1,1-dimethylethyl)dimethylsilyl]oxy]-,
 phenylmethyl ester (9CI)

MF C21 H28 O3 Si

$$\begin{array}{c|c} O & \\ \parallel & \\ CH_2-C-O-CH_2-Ph \\ \downarrow & \\ t-Bu-Si-O & \\ Me & \\ Me & \\ \end{array}$$

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L1 9 ANSWERS REGISTRY COPYRIGHT 2007 ACS on STN

IN Benzenepropanoic acid, β-(aminocarbonyl)-4-bromo- (9CI)

MF C10 H10 Br N O3

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L1 9 ANSWERS REGISTRY COPYRIGHT 2007 ACS on STN

IN Benzeneacetic acid, α -(2-amino-2-oxoethyl)-4-bromo- (9CI)

MF C10 H10 Br N O3

Page 4 searched 9/5/07 updated str search

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L1 9 ANSWERS REGISTRY COPYRIGHT 2007 ACS on STN

IN Benzenepropanoic acid, β -(aminocarbonyl)-4-bromo-, phenylmethyl ester (9CI)

MF C17 H16 Br N O3

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L1 9 ANSWERS REGISTRY COPYRIGHT 2007 ACS on STN

IN Benzeneacetic acid, α -(2-amino-2-oxoethyl)-4-bromo-, phenylmethyl

ester (9CI)

MF C17 H16 Br N O3

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L1 9 ANSWERS REGISTRY COPYRIGHT 2007 ACS on STN

IN Butanedioic acid, (4-bromophenyl)-, 4-(phenylmethyl) ester (9CI)

MF C17 H15 Br O4

Page 5 searched 9/5/07 updated str search

ALL ANSWERS HAVE BEEN SCANNED

=> d his

(FILE 'HOME' ENTERED AT 18:07:18 ON 25 MAR 2007)

FILE 'REGISTRY' ENTERED AT 18:07:32 ON 25 MAR 2007

L1 9 S 335200-36-7/RN OR 845785-97-9/RN OR 845785-98-0/RN OR 8457

=> fil hcap

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION

FULL ESTIMATED COST 0.45 0.66

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=> s 11

L2 2 L1

=> d 12 1-2 ibib abs

L2 ANSWER 1 OF 2 HCAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER: 2005:158625 HCAPLUS

DOCUMENT NUMBER:

142:261292

TITLE:

LANGUAGE:

Preparation of (hetero)aryl-substituted succinate derivatives as matrix metalloproteinase inhibitors

INVENTOR(S): Holmes, Ian; Watson, Stephen Paul

PATENT ASSIGNEE(S): Glaxo Group Limited, UK SOURCE: PCT Int. Appl., 36 pp.

CODEN: PIXXD2

DOCUMENT TYPE:

Patent English

FAMILY ACC. NUM. COUNT: 1

Page 6 searched 9/5/07 updated str search

PATENT INFORMATION:

PATENT NO.					KIND		DATE			APPL	ICAT	ION	DATE					
_	2005016868 2005016868								WO 2004-EP9087						20040812			
	W:						AU,											
		CN,	CO,	CR,	CU,	CZ,	DE,	DK,	DM,	DZ,	EC,	EE,	EG,	ES,	FI,	GB,	GD,	
		GE,	GH,	GM,	HR,	HU,	ID,	ΙL,	IN,	IS,	JP,	KE,	KG,	KP,	KR,	KZ,	LC,	
		LK,	LR,	LS,	LT,	LU,	LV,	MA,	MD,	MG,	MK,	MN,	MW,	MX,	MZ,	NA,	NI,	
							PL,											
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	RW:						MW,											
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EP	EP 1654218				A2 20060510			EP 2004-764084					20040812					
		_		СH			ES,											
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at.	IE, SI, LT, JP 2007502259																	
US 2006235074																		
PRIORITY APPLN. INFO.:				A1 20061019				GB 2003-19069										
FRIORITI AFPLIN. INFO.:																		
OMITTED COLLEGE (C)					WO 2004-EP9087 W 20040812									512				
					CASREACT 142:261292; MARPAT 142:261292													
GI																		

Title compds. represented by the formula I, R1ZQCH(R2)CH2X, [wherein R1 = (un)substituted alkyl(cycloalkyl), alkylheterocycloalkyl, alkylaryl, etc.; Z = a bond, CH2, O, S, etc.; Q = (un)substituted (hetero)aryl; X = COR3; R2 = CONH2, CO2H, sulfonylamino, etc.; R3 = OH, oxyalkyl or (un)substituted amino; with a proviso; and physiol. functional derivs. thereof] were prepared as matrix metalloproteinase (MMP) inhibitors. Coupling reaction of 4-amino-3-(4-bromophenyl)-4-oxobutanoic acid with p-nitrilephenylboronic acid gave II in 100% yield. I showed inhibition of MMP-12 with IC50 values of below 100 μ M. Thus, I and their pharmaceutical compns. are useful as matrix metalloproteinase inhibitors for the treatment of inflammation or autoimmune disease (no data).

L2 ANSWER 2 OF 2 HCAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER: 2001:8060 HCAPLUS

DOCUMENT NUMBER: 134:307022

TITLE: Antibody-catalyzed hydrolysis of oligomeric esters: a

model for the degradation of polymeric materials

AUTHOR(S): Brummer, Oliver; Hoffman, Timothy Z.; Chen, Da-Wei;

Janda, Kim D.

CORPORATE SOURCE: Department of Chemistry, The Scripps Research

Institute and The Skaggs Institute for Chemical

Biology, La Jolla, CA, 92037, USA

SOURCE: Chemical Communications (Cambridge) (2001), (1), 19-20

CODEN: CHCOFS; ISSN: 1359-7345

PUBLISHER:

```
DOCUMENT TYPE:
                         Journal
LANGUAGE:
                         English
OTHER SOURCE(S):
                         CASREACT 134:307022
    A catalytic antibody has been discovered that degrades oligomeric ester
     substrates. All the observations and data confirmed that the antibody
     performed oligomer degrdns. by 'multimer' processing using
     nonregioselective, kinetically biased endo-cleavage, rather than a
     stepwise deoligomerization through cleavage of monomers from a terminus.
     These findings are of fundamental importance as now catalytic antibodies
     share another trait thought only to be associated with enzymes, the
     biodegrdn. of oligo and polymeric materials.
REFERENCE COUNT:
                         14
                               THERE ARE 14 CITED REFERENCES AVAILABLE FOR THIS
                               RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT
=> s 845786-06-3/rn or 845786-07-4/rn or 845786-08-5/rn or 845786-09-6/rn or
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               OR 845786-13-2/RN OR 845786-14-3/RN
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Royal Society of Chemistry

=> file reg

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION 11.52

10.86 FULL ESTIMATED COST

TOTAL DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE ENTRY SESSION

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23 MAR 2007 HIGHEST RN 928114-47-0 STRUCTURE FILE UPDATES: DICTIONARY FILE UPDATES: 23 MAR 2007 HIGHEST RN 928114-47-0

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http://www.cas.org/ONLINE/UG/regprops.html

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845786-06-3 IS NOT A RECOGNIZED COMMAND

The previous command name entered was not recognized by the system. For a list of commands available to you in the current file, enter "HELP COMMANDS" at an arrow prompt (=>).

=> s 845786-06-3/rn or 845786-07-4/rn or 845786-08-5/rn or 845786-09-6/rn or 845786-10-9/rn or 845786-11-0/rn or 845786-12-1/rn or 845786-13-2/rn or 845786-14-3/rn

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1 845786-13-2/RN

1 845786-14-3/RN

9 845786-06-3/RN OR 845786-07-4/RN OR 845786-08-5/RN OR 845786-L4 09-6/RN OR 845786-10-9/RN OR 845786-11-0/RN OR 845786-12-1/RN OR 845786-13-2/RN OR 845786-14-3/RN

=> d scan

L4 9 ANSWERS REGISTRY COPYRIGHT 2007 ACS on STN

IN Benzeneacetic acid, α -(2-amino-2-oxoethyl)-4-hydroxy- (9CI)

MF C10 H11 N O4

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):9

L4 9 ANSWERS REGISTRY COPYRIGHT 2007 ACS on STN

IN Butanedioic acid, [4-[[(1,1-dimethylethyl)dimethylsilyl]oxy]phenyl]-,

1-(phenylmethyl) ester (9CI)

MF C23 H30 O5 Si

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L4 9 ANSWERS REGISTRY COPYRIGHT 2007 ACS on STN

IN Butanedioic acid, [4-(3-methylbutoxy)phenyl]-, 4-(1,1-dimethylethyl)

1-methyl ester (9CI)

MF C20 H30 O5

Page 10 searched 9/5/07 updated str search

L4 9 ANSWERS REGISTRY COPYRIGHT 2007 ACS on STN

MF C22 H26 O5

$$\begin{array}{c|c} & CO_2H & O \\ & | & | \\ CH-CH_2-C-O-CH_2-Ph \\ \\ \text{Me}_2\text{CH}-CH_2-CH_2-O \end{array}$$

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L4 9 ANSWERS REGISTRY COPYRIGHT 2007 ACS on STN

IN Benzeneacetic acid, 4-(3-methylbutoxy)-, 1,1-dimethylethyl ester (9CI)

MF C17 H26 O3

$$\begin{array}{c} \text{O} \\ \text{CH}_2-\text{C-OBu-t} \\ \text{Me}_2\text{CH}-\text{CH}_2-\text{CH}_2-\text{O} \end{array}$$

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L4 9 ANSWERS REGISTRY COPYRIGHT 2007 ACS on STN

IN Benzeneacetic acid, α -(2-amino-2-oxoethyl)-4-hydroxy-, phenylmethyl

ester (9CI)

MF C17 H17 N O4

Page 11 searched 9/5/07 updated str search

L4 9 ANSWERS REGISTRY COPYRIGHT 2007 ACS on STN

IN Butanedioic acid, [4-[[(1,1-dimethylethyl)dimethylsilyl]oxy]phenyl]-,

4-(1,1-dimethylethyl) 1-(phenylmethyl) ester (9CI)

MF C27 H38 O5 Si

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L4 9 ANSWERS REGISTRY COPYRIGHT 2007 ACS on STN

IN Benzeneacetic acid, 4-(3-methylbutoxy)-, methyl ester (9CI)

MF C14 H20 O3

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L4 9 ANSWERS REGISTRY COPYRIGHT 2007 ACS on STN

IN Butanedioic acid, [4-(3-methylbutoxy)phenyl]-, 1-(1,1-dimethylethyl)

4-(phenylmethyl) ester (9CI)

MF C26 H34 O5

$$\begin{array}{c|c} & & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ \text{Me}_2\text{CH}-\text{CH}_2-\text{CH}_2-\text{O} \end{array}$$

Page 12 searched 9/5/07 updated str search

ALL ANSWERS HAVE BEEN SCANNED

=> fil hcap COST IN U.S. DOLLARS SINCE FILE TOTAL SESSION ENTRY 11.97 FULL ESTIMATED COST 0.45 TOTAL SINCE FILE DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) ENTRY SESSION 0.00 CA SUBSCRIBER PRICE -1.56

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FILE COVERS 1907 - 25 Mar 2007 VOL 146 ISS 14 FILE LAST UPDATED: 23 Mar 2007 (20070323/ED)

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This file contains CAS Registry Numbers for easy and accurate substance identification.

=> d his

(FILE 'HOME' ENTERED AT 18:07:18 ON 25 MAR 2007)

FILE 'REGISTRY' ENTERED AT 18:07:32 ON 25 MAR 2007

L1 9 S 335200-36-7/RN OR 845785-97-9/RN OR 845785-98-0/RN OR 8457

FILE 'HCAPLUS' ENTERED AT 18:08:19 ON 25 MAR 2007

L2 2 S L1

L3 1 S 845786-06-3/RN OR 845786-07-4/RN OR 845786-08-5/RN OR 8457

FILE 'REGISTRY' ENTERED AT 18:09:41 ON 25 MAR 2007

L4 9 S 845786-06-3/RN OR 845786-07-4/RN OR 845786-08-5/RN OR 8457

FILE 'HCAPLUS' ENTERED AT 18:10:04 ON 25 MAR 2007

=> s 14

L5 1 L4

Page 13 searched 9/5/07 updated str search

=> d ibib abs

ANSWER 1 OF 1 HCAPLUS COPYRIGHT 2007 ACS on STN

2005:158625 HCAPLUS ACCESSION NUMBER:

142:261292 DOCUMENT NUMBER:

Preparation of (hetero)aryl-substituted succinate TITLE:

derivatives as matrix metalloproteinase inhibitors

Holmes, Ian; Watson, Stephen Paul Glaxo Group Limited, UK INVENTOR (S):

PATENT ASSIGNEE(S): PCT Int. Appl., 36 pp. SOURCE:

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PATENT				TE APPLICATION NO.							DATE					
					WO 2004 ED0007						20040812					
									WO 2004-EP9087							
WO 2005	A3		2005	0519												
W:	AE, AG,	AL,	AM,	AT,	AU,	ΑZ,	BA,	BB,	BG,	BR,	BW,	BY,	ΒZ,	CA,	CH,	
	CN, CO,	CR,	CU,	CZ,	DE,	DK,	DM,	DZ,	EC,	EE,	EG,	ES,	FI,	GB,	GD,	
	GE, GH,	GM,	HR,	HU,	ID,	IL,	IN,	IS,	JΡ,	KE,	KG,	KP,	KR,	ΚZ,	LC,	
	LK, LR,															
	NO, NZ,															
	TJ, TM,															
RW:	BW, GH,															
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	EE, ES,															
	SI, SK,															
	SN, TD,	TG														
EP 1654	A2 20060510]	EP 2	004-	20040812								
R:	AT, BE,	CH,	DE,	DK,	ES,	FR,	GB,	GR,	IT,	LI,	LU,	NL,	SE,	MC,	PT,	
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JP 200						0208	JP 2006-522996						20040812			
US 2006	A1		2006	1019	1	US 2	006-	5698	12		2	0060	210			
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							1	WO 2	004-	EP90	87	1	W .2	0040	812	
OTHER SOURCE	CASREACT 142:261292; MARPAT 142:261292															
GI																

Title compds. represented by the formula I, R1ZQCH(R2)CH2X, [wherein R1 = AΒ (un)substituted alkyl(cycloalkyl), alkylheterocycloalkyl, alkylaryl, etc.; Z = a bond, CH2, O, S, etc.; Q = (un)substituted (hetero)aryl; X = COR3; R2 = CONH2, CO2H, sulfonylamino, etc.; R3 = OH, oxyalkyl or (un) substituted amino; with a proviso; and physiol. functional derivs. thereof] were prepared as matrix metalloproteinase (MMP) inhibitors. Coupling reaction of 4-amino-3-(4-bromophenyl)-4-oxobutanoic acid with

Page 14 searched 9/5/07 updated str search

p-nitrilephenylboronic acid gave II in 100% yield. I showed inhibition of MMP-12 with IC50 values of below 100 μM . Thus, I and their pharmaceutical compns. are useful as matrix metalloproteinase inhibitors for the treatment of inflammation or autoimmune disease (no data).

=> fil stng SINCE FILE COST IN U.S. DOLLARS TOTAL ENTRY SESSION FULL ESTIMATED COST 10.63 22.60 DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL SESSION ENTRY -0.78 -2.34 CA SUBSCRIBER PRICE

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